

Application No.: 10/517713

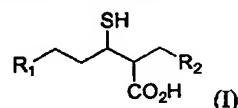
Docket No.: ASZD-P01-722

Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

1. (currently amended) A compound of formula (I):



wherein:

R¹ is phenyl {optionally substituted by halogen, hydroxy, cyano, C₁₋₄ alkyl (itself optionally mono-substituted by cyano, hydroxy or phenyl), C₁₋₄ alkoxy (itself optionally substituted by tetrahydrofuranyl), CF₃, OCF₃, methylenedioxy, C(O)R³, S(O)₂R⁴, phenyl (itself optionally substituted by halogen), phenoxy (itself optionally substituted by halogen) or tetrahydrofuryloxy}, naphthyl, pyridinyl, 1,2,3,4-tetrahydropyrimidin-2,4-dione-yl (optionally substituted by C₁₋₄ alkyl) or tetrahydrothienyl;

R² is aminopyridinyl, aminothiazolyl or 3-azabicyclo[3.2.1]octyl;

R³ is hydroxy, C₁₋₄ alkoxy (itself optionally substituted by phenyl (itself optionally substituted by halogen) or pyridinyl), NR⁵R⁶ or an N-linked 5- or 6-membered heterocyclic ring {unsubstituted or mono-substituted by hydroxy, oxo, C₁₋₄ alkyl (itself optionally substituted by hydroxy or NHphenyl), CO²(C₁₋₄ alkyl) or phenyl (itself optionally substituted by halogen)};

R⁴ is NR⁷R⁸ or an N-linked 5- or 6-membered heterocyclic ring {unsubstituted; mono-substituted by hydroxy, oxo, C₁₋₄ alkyl (itself optionally substituted by hydroxy or NHphenyl), CO₂(C₁₋₄ alkyl) or phenyl (itself optionally substituted by halogen); or fused to a benzene ring which is optionally substituted by C₁₋₄ alkoxy};

R⁵, R⁶, R⁷ and R⁸ are, independently, hydrogen, C₁₋₄ alkyl {optionally substituted by halogen, cyano, hydroxy, phenyl (itself optionally substituted by halogen or methylenedioxy), pyridinyl, CO₂H or CO₂(C₁₋₄ alkyl)} I or C₂₋₄ alkenyl;

Application No.: 10/517713

Docket No.: ASZD-P01-722

provided that when \mathbf{R}^1 \mathbf{R}^2 is 6-aminopyridin-3-yl then \mathbf{R}^2 \mathbf{R}^1 is substituted phenyl, naphthyl, pyridinyl, 1,2,3,4-tetrahydropyrimidin-2,4-dione-yl (optionally substituted by C_{1-4} alkyl) or tetrahydrothienyl; or a pharmaceutically acceptable salt or solvate thereof, or a solvate of such a salt.

2. (original) A compound of formula (I) as claimed in claim 1 wherein \mathbf{R}^1 is phenyl (optionally substituted by halogen, hydroxy, cyano, C_{1-4} alkyl (itself optionally mono-substituted by cyano or hydroxy), C_{1-4} alkoxy, CF_3 , OCF_3 , methylenedioxy, $\text{C}(\text{O})\text{NH}_2$, $\text{S}(\text{O})_2\text{NH}_2$ or phenyl (itself optionally substituted by halogen)), pyridinyl or tetrahydrothienyl.

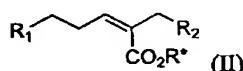
3. (original) A compound of formula (I) as claimed in claim 1 wherein \mathbf{R}^1 is phenyl (optionally substituted by halogen, hydroxy, cyano, C_{1-4} alkyl (itself optionally mono-substituted by cyano, hydroxy or phenyl), C_{1-4} alkoxy, CF_3 , OCF_3 , methylenedioxy, phenoxy (itself optionally substituted by halogen), tetrahydrofuryloxy or tetrahydrofuranmethoxy), naphthyl, pyridinyl or tetrahydrothienyl.

4. (original) A compound of formula (I) as claimed in claim 1 wherein \mathbf{R}^1 is phenyl (substituted by halogen, hydroxy, cyano, C_{1-4} alkyl (itself optionally mono-substituted by cyano or hydroxy), C_{1-4} alkoxy, CF_3 or methylenedioxy) or tetrahydrothiophenyl.

5. (original) A compound of formula (I) as claimed in claim 1, 2, 3 or 4 wherein \mathbf{R}^2 is 6-aminopyridin-3-yl, 2-aminothiazol-5-yl or 3-azabicyclo[3.2.1]oct-8-yl.

6. (original) A compound of formula (I) as claimed in claim 1, 2, 3 or 4 wherein \mathbf{R}^2 is 6-aminopyridin-3-yl.

7. (original) A process for preparing a compound of formula (I) comprising reacting a compound of formula (II):

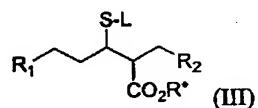


wherein \mathbf{R}^1 is as defined in claim 1 or includes a group that can be subsequently reacted

Application No.: 10/517713

Docket No.: ASZD-P01-722

to form the group $R^1 R^*$ is a suitable protecting group and R^2 is as defined in claim 1 or the amine function of R^2 can be protected, with a thiol of formula $L-SH$, wherein L is a suitable protecting group, in the presence of a suitable catalyst and in a suitable solvent, to form a compound of formula (III):



and, optionally reacting the functional group on R^1 , and subsequently removing the protecting groups as necessary.

8. (original) A pharmaceutical formulation containing a compound according to any one of claims 1 to 6 as active ingredient in combination with a pharmaceutically acceptable adjuvant, diluent or carrier.

9. (original) The use of a compound as claimed in claim 1 in therapy.

10. (original) The use of a compound as claimed in claim 1 for the manufacture of a medicament for the inhibition of carboxypeptidase U.

11. (original) A method for treatment or prophylaxis of conditions where inhibition of carboxypeptidase U is beneficial, comprising administering to a mammal, including man, in need of such treatment an effective amount of a compound as claimed in claim 1.

12. (original) A pharmaceutical formulation for use in the treatment or prophylaxis of conditions where inhibition of carboxypeptidase U is beneficial, comprising a compound as claimed in claim 1 in combination with a pharmaceutically acceptable adjuvant, diluent or carrier.